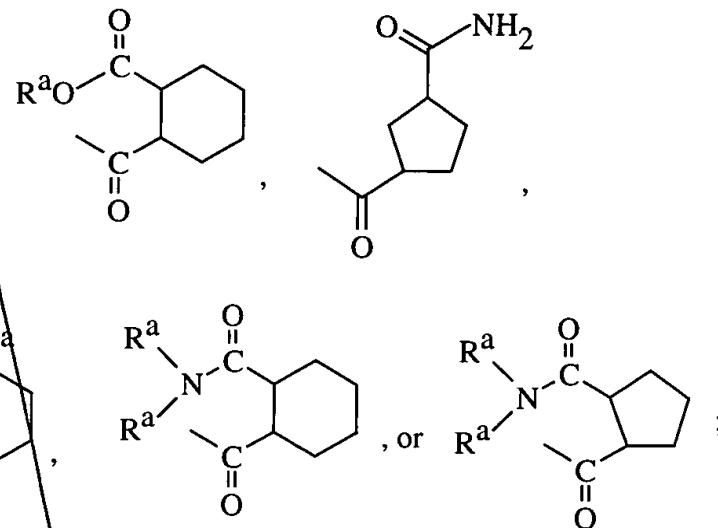


$R^3CO^-$ ,  
 $R^3SO_2^-$ ,

$R^a$

$R^5NCHR^6CO^-$ ,



each  $R^a$  is independently hydrogen,  $C_1-C_6$  alkyl, or  $-(CH_2)_n$  aryl;

$R^2$  is  $-(CRR)_n$ -aryl,

$-(CRR)_n$ -X-aryl,

$-(CRR)_n$ -heteroaryl,

$-(CRR)_n$ -X-heteroaryl,

$-(CRR)_n$ -(substituted-heteroaryl),

$-(CRR)_n$ -(substituted-aryl),

$-(CRR)_n$ -X-(substituted-aryl),

$-(CRR)_n$ -aryl-aryl,

$-(CRR)_n$ -aryl-heteroaryl,

$-(CRR)_n$ -aryl- $(CH_2)_n$ -aryl,

$-(CRR)_n$ -CH(aryl) $_2$ ,

$-(CRR)_n$ -cycloalkyl,

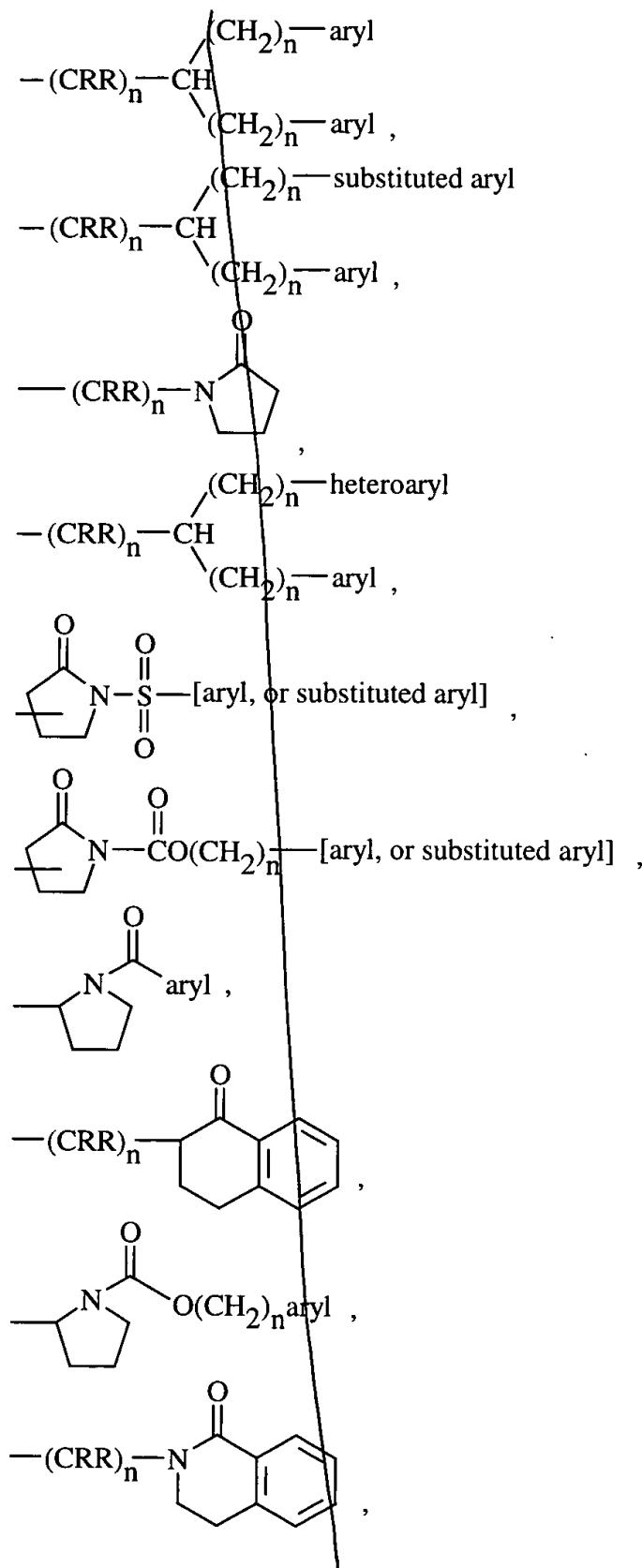
$-(CRR)_n$ -X-cycloalkyl,

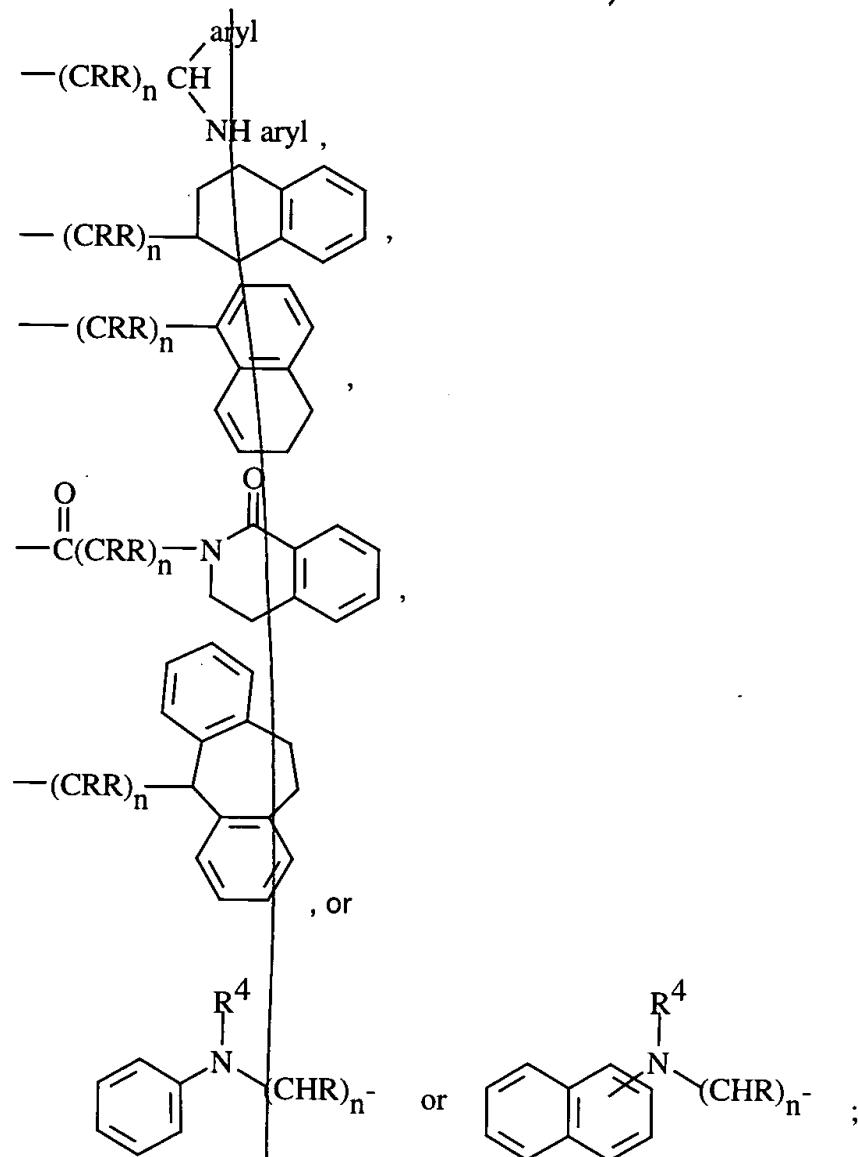
$-(CRR)_n$ -heterocycle,

$-(CRR)_n$ -X-heterocycle,

$-(CRR)_n$  substituted heterocycle,

B  
a  
cont'd





each R is independently hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, halogen or hydroxy;

X is O or S;

R<sup>3</sup> is C<sub>1</sub>-C<sub>6</sub> alkyl,

aryl,

heteroaryl,

-(CHR)<sub>n</sub>-aryl,

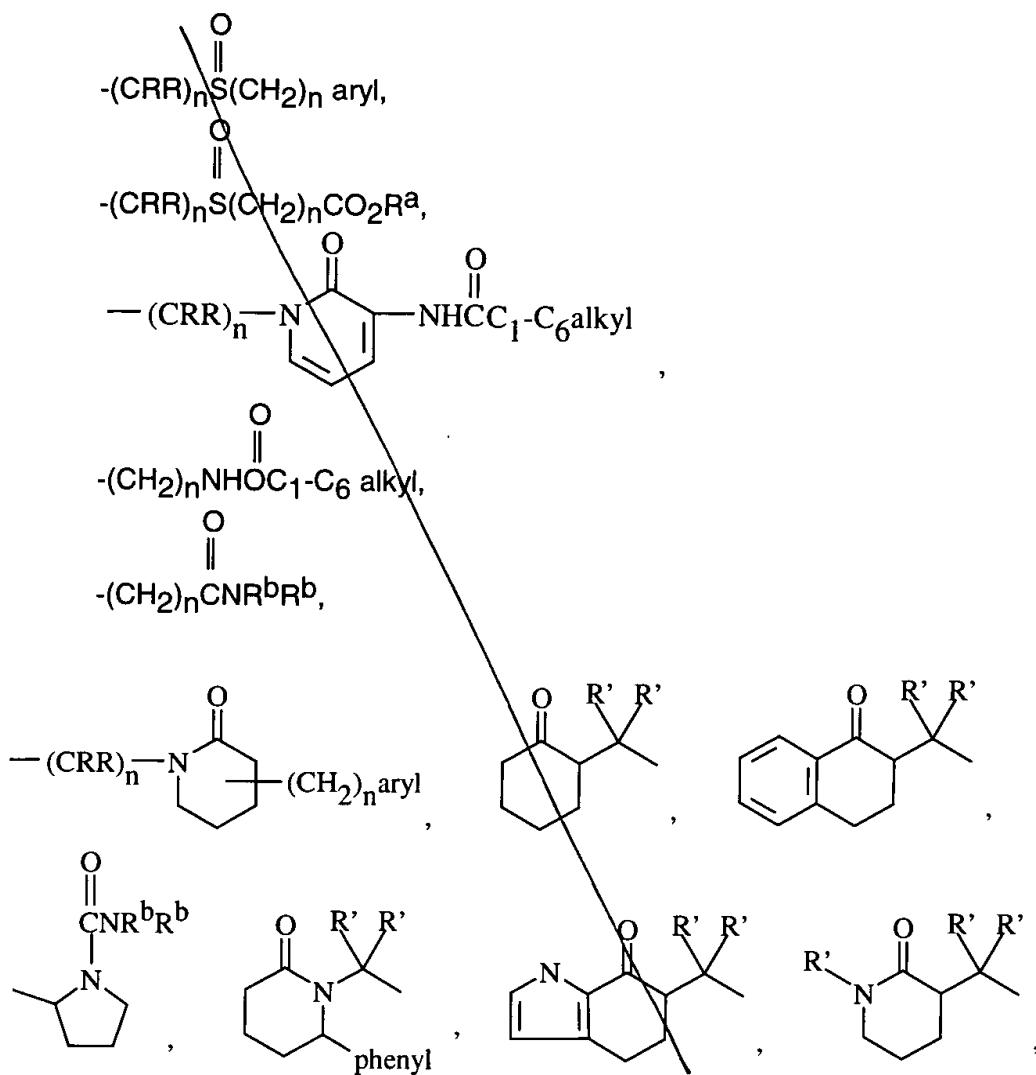
-(CHR)<sub>n</sub>-heteroaryl,

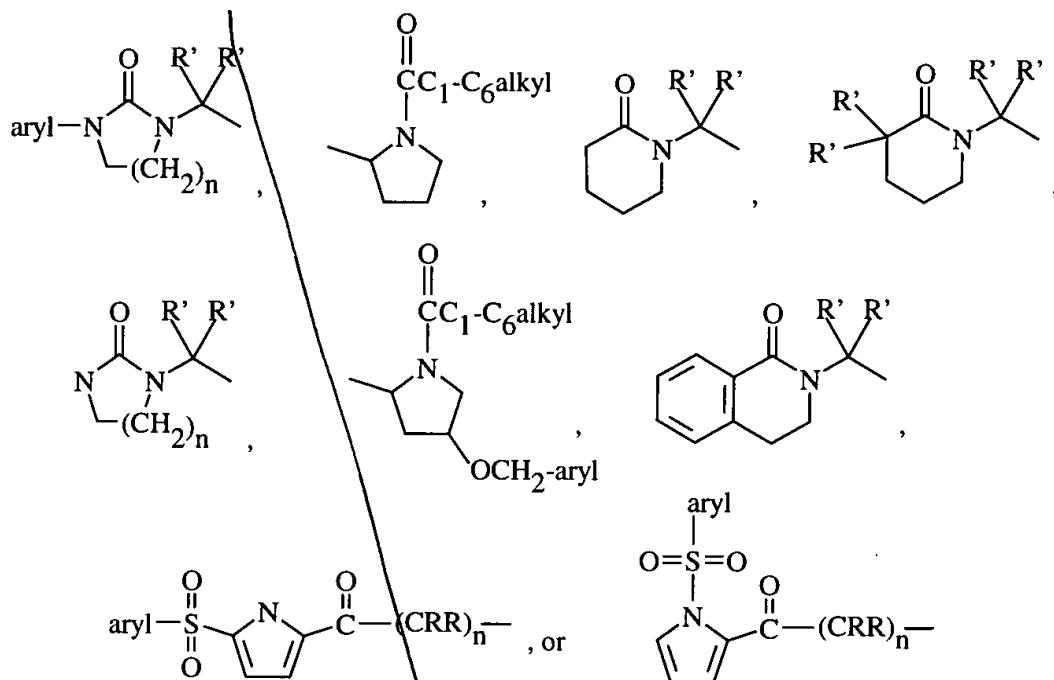
-(CHR)<sub>n</sub>-substituted heteroaryl,

-(CHR)<sub>n</sub>-substituted aryl,

$-(CRR)_nCOR^a$ ,  
 $-(CRR)_nS(CH_2)_n$ -aryl,  
cycloalkyl,  
substituted cycloalkyl,  
heterocycle,  
substituted heterocycle,  
 $-(CRR)_nCNR^aR^a$ ,  
 $-(CRR)_n-S-(CH_2)_n$  aryl,  
 $-(CRR)_n-SC_1-C_6$  alkyl,  
 $J-CH_2CH-$ ,  
 $-CHR^6C$ -heteraryl,  
 $-(CRR)_nS(CH_2)_nCOR^a$ ,  
 $-(CRR)_nS(CH_2)_nCOR^a$ ,  
 $-(CRR)_nS(CH_2)_n$ -aryl,  
 $-(CRR)_nS(CH_2)_n$ -aryl,  
 $-(CRR)_nSCC_1-C_6$  alkyl,

*✓  
contd*





each R' is independently C<sub>1</sub>-C<sub>6</sub> alkyl,

C<sub>1</sub>-C<sub>6</sub> alkylaryl,

aryl, or

hydrogen;

each J is independently

-CO<sub>2</sub>R<sup>b</sup>,

-CONR<sup>b</sup>R<sup>b</sup>,

-SO<sub>2</sub>NR<sup>b</sup>R<sup>b</sup>, or

-SO<sub>2</sub>R<sup>b</sup>;

each R<sup>b</sup> is independently hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, aryl, substituted aryl; arylalkyl,

heteroarylalkyl, substituted arylalkyl, or substituted heteroarylalkyl;

R<sup>4</sup> is hydrogen,

C<sub>1</sub>-C<sub>6</sub> alkyl,

CH<sub>3</sub>OC-,

-phenyl, or

*Art  
contd*

*R<sup>5</sup> is*

$\text{C}_1\text{-C}_6$  alkyl C-;

$\text{C}_1\text{-C}_6$  alkyl-CO-;

$-(\text{CH}_2)_n$  aryl,

$\text{C}_1\text{-C}_6$ -alkyLOC-;

$\text{C}_1\text{-C}_6$ -alkyl-X-(CH<sub>2</sub>)<sub>n</sub>CO,

$\text{C}_1\text{-C}_6$ -alkyl-X-(CH<sub>2</sub>)<sub>n</sub>OC-;

$-\text{C}(\text{CRR})_n$  aryl,

$-\text{CNR}^a\text{R}^a$ ,

$-\text{SC}_1\text{-C}_6$  alkyl,

$-\text{C}(\text{CH}_2)_n\text{CNR}^a\text{R}^a$ ,

$-\text{CO}(\text{CH}_2)_n$  aryl,

$-\text{CO}(\text{CH}_2)_n$  substituted aryl,

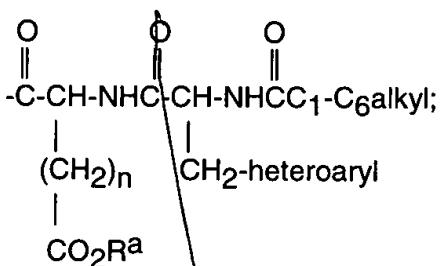
$-\text{C}(\text{CRR})_n\text{NHCO}(\text{CH}_2)_n$ -aryl,

$\text{C}(\text{R}^6)-\text{CH}-\text{N}(\text{R}^a)\text{R}^{5a}$

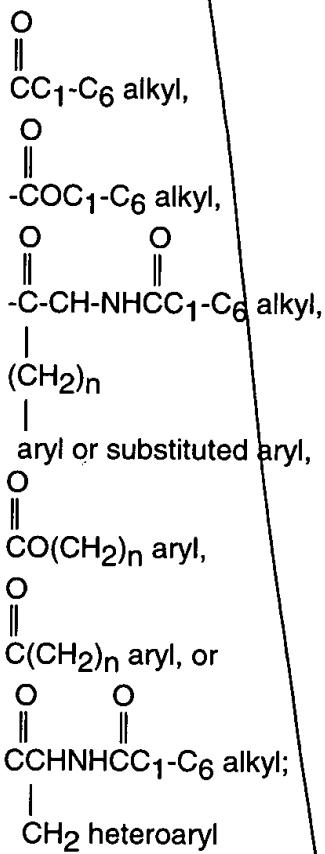
$-(\text{CH}_2)_n\text{X}(\text{CH}_2)_n$ -aryl,

$-\text{C}_1\text{-C}_6$  alkyl X-C<sub>1</sub>-C<sub>6</sub> alkyl aryl, or

*B*  
*at*  
*Contd*



$\text{R}^5\text{a}$  is

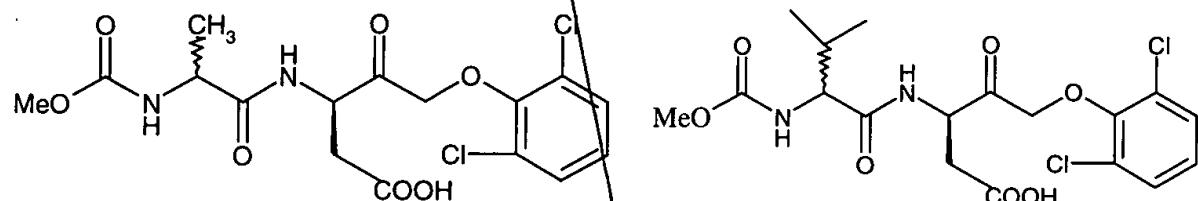


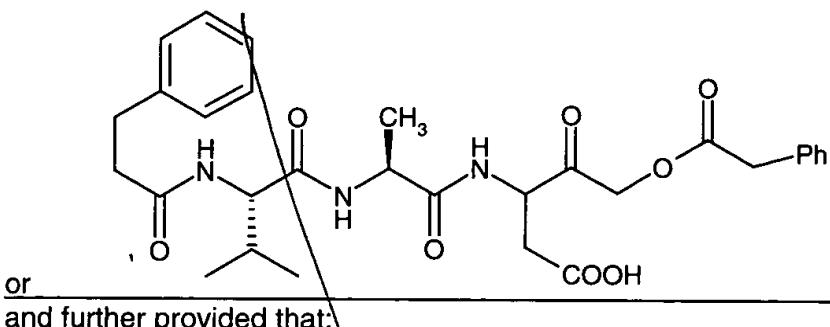
$\text{R}^6$  is hydrogen,

$\text{C}_1\text{-C}_6$  alkyl,  $-(\text{CH}_2)_n$  aryl,  $-(\text{CH}_2)_n\text{CO}_2\text{R}^a$ , hydroxyl substituted  $\text{C}_1\text{-C}_6$  alkyl, or imidazole substituted  $\text{C}_1\text{-C}_6$  alkyl;

each  $n$  is independently 0 to 3, and the pharmaceutically acceptable, salts, esters, amides, and prodrugs thereof;

provided that a compound of formula (I) is not:





or

and further provided that:

(a) when  $R^2$  is aryl, substituted aryl, cycloalkyl, phenyl-phenyl- $CH_2$ -, piperidino, heteroaryl or substituted heteroaryl; and  $R^1$  is  $(R^5R^a)N-CH(R^6)-CO-$ , then  $R^a$  is not hydrogen when

$R^6$  is a side chain of an amino acid;

$R^5$  is aryl- $C(O)-$ , aryl- $(CH_2)-O-C(O)-NH-CH(R)-C(O)-$ ;

where  $R$  is H or  $(C_1-C_6)$ alkyl or  $R^{5a}-NH-CH(R^6)-C(O)-$ ;

where  $R^6$  is a side chain of an amino acid and  $R^{5a}$  is an amino acid protecting group;

(b) when  $R^1$  is  $R^3-O-C(O)-$  where  $R^3$  is  $CH_2=CH-CH_2$ -, then  $R^2$  is not  $Ph(CH_2)_2$ -,  $PhO(CH_2)_2$ -, trans- $PhCH=CH$  or cyclohexyl( $CH_2)_2$ ;

(c) when  $R^1$  is  $(R^5R^a)N-CH(R^6)-CO-$ ;

where  $R^6$  is H,  $(C_1-C_6)$ alkyl, benzyl or hydroxyalkyl;

$R^a$  is H,  $(C_1-C_6)$ alkyl, phenyl or benzyl; and

$R^5$  is  $-C(O)-O-(C_1-C_6)$ alkyl,  $-C(O)-N(R^aR^a)$ ,  $-C(O)-(C^1-C^6)$ alkyl, -phenyl- $O-(C_1-C_6)$ alkyl or -phenyl- $(CH_2)_{1-4}-N(R^aR^a)$ ;

then  $R^2$  is not a phenyl or naphthyl group optionally substituted with one or more substituents selected from the group consisting of halogen, hydroxy,  $CF_3$ ,  $NO_2$ ,  $(C_1-C_6)$ alkoxy,  $-CO-(C_1-C_6)$ alkyl,  $-NR^aC(O)-(C_1-C_6)$ alkyl,  $-CON(R^aR^a)$ ,  $-SO_2N(R^aR^a)$ ,  $-SO_2-(C_1-C_6)$ alkyl,  $-COO-(C_1-C_6)$ alkyl,  $(C_1-C_6)$ alkyl, cycloalkyl and  $-O-(CH_2)_{1-6}-phenyl-O-(C_1-C_6)$ alkyl; and

(d) when  $R^1$  is  $R^5-NH-CH(R^6)-C(O)-$ , where  $R^5$  is  $R^{5a}-NH-CH(R^6)-C(O)-$  and  $R^{5a}$  is  $-C(O)-(C_1-C_6)$ alkyl or  $-C(O)-aryl$ , then  $R^2$  is not mono-, di-, tri-, tetra- or penta-substituted phenyl or mono-, di-, tri-substituted phenyl, 1-naphthyl, 9-anthracyl or 2-, 3- or 4-pyridyl.

Please cancel claim 9, without waiver or prejudice.

Please cancel claim 32, without waiver or prejudice, and insert therefor new claim 44:

--44. The compounds:

3-Benzoyloxycarbonylamino-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-Benzoyloxycarbonylamino-4-oxo-5-(3-phenyl-propionyloxy)-pentanoic acid;

3-Benzylcarbonylamino-5-(3-cyclohexyl-propionyloxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-[(naphthalene-1-yl-oxy)-acetoxy]-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-4-oxo-5-phenoxyacetoxy-pentanoic acid;  
3-Benzylcarbonylamino-4-oxo-5-phenylsulfanylacetoxypentanoic acid;  
3-Benzylcarbonylamino-5-[(6-methoxy-naphthalene-1-yl)-acetoxy]-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(naphthalene-2-yl-acetoxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(3-naphthalene-2-yl-propionyloxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(3,3-diphenyl-propionyloxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(2-naphthalene-1-yl-propionyloxy)-4-oxo-pentanoic acid;  
5-[(Acetyl-phenyl-amino)-acetoxy]-3-benzylcarbonyl-amino-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-(hydroxy-naphthalene-1-yl-acetoxy)-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-4-oxo-5-[(phenyl-amino)-acetoxy]-pentanoic acid;  
3-Benzylcarbonylamino-5-[(6-hydroxy-naphthalene-1-yl)-acetoxy]-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-5-[3-(4-hydroxy-phenyl)-2-naphthalene-1-yl-propionyloxy]-4-oxo-pentanoic acid;  
(S)-3-Benzylcarbonylamino-4-oxo-5-phenylacetoxypentanoic acid;  
(S)-3-Benzylcarbonylamino-4-oxo-5-(4-phenyl-butyryloxy)-pentanoic acid;  
3-Benzylcarbonylamino-4-oxo-5-[(4-phenyl-naphthalen-1-yl)-acetoxy]-pentanoic acid;  
3-Benzylcarbonylamino-5-[(4-methyl-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;  
3-Benzylcarbonylamino-4-oxo-5-[(4-thiophen-2-yl-naphthalen-1-yl)-acetoxy]-pentanoic acid;  
3-Benzylcarbonylamino-5-[(4-fluoro-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

*CH 1  
CONT'D*

3-Benzylloxycarbonylamino-5-[(2-methyl-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(2-fluoro-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

5-[(4-Benzyl-naphthalen-1-yl)-acetoxy]-3-benzylloxycarbonylamino-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(3,4-dihydro-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-(3,4-diphenyl-butyryloxy)-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-4-oxo-5-(3-phenyl-3-phenylamino-propionyloxy)-pentanoic acid;

3-Benzylloxycarbonylamino-4-oxo-5-[(1,2,3,4-tetrahydro-naphthalen-2-yl)-acetoxy]-pentanoic acid;

3-Benzylloxycarbonylamino-4-oxo-5-[(2,3,5,6-tetramethyl-phenyl)-acetoxy]-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(2,3-dichloro-phenyl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(5-methyl-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(2-iodo-phenyl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(5-methoxy-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(8-methyl-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(9H-fluoren-9-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-[(10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-yl)-acetoxy]-4-oxo-pentanoic acid;

3-Benzylloxycarbonylamino-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid; and

3-Benzylloxycarbonylamino-5-[(5-cyano-naphthalen-1-yl)-acetoxy]-4-oxo-pentanoic acid.--

Please cancel claim 33, without waiver or prejudice and insert therefor new claim 45:

45. The compounds:

[S-(R\*,R\*)]-3-(2-Acetylamino-propionylamino)-5-(naphthalene-1-yl-acetoxy)-4-oxo-pentanoic acid;

5-(Naphthalen-1-yl-acetoxy)-4-oxo-3-[2-(4-phenyl-butyrylamino)-propylamino]-pentanoic acid;

3-(2-Methanesulfonylamino-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-[2-(2-Acetylamino-4-phenyl-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Acetylamino-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-[2-(4-Carbamoyl-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Benzylloxycarbonylamino-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

5-(Naphthalen-1-yl-acetoxy)-4-oxo-3-(2-ureido-propionylamino)-pentanoic acid;

3-(2-Acetylamino-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Acetylamino-acetylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Acetylamino-propionylamino)-5-(3,3-diphenyl-propionyloxy)-4-oxo-pentanoic acid;

3-[2-(2-Acetylamino-4-carboxy-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

5-(Naphthalen-1-yl-acetoxy)-4-oxo-3-[2-(3-phenyl-propionylamino)-propionylamino]-pentanoic acid;

3-[2-(3-Methyl-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid; and

3-(4-Carbamoyl-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid.--

Please cancel claim 35, without waiver or prejudice, and insert therefor new claim 46:

--46. The compounds:

3-(2-Methyl-3-phenethylcarbamoyl-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-[3-Methyl-2-(3-phenyl-propionylamino)-butyrylamino]-4-oxo-5-[(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetoxy]-pentanoic acid;  
5-(Naphthalen-1-yl-acetoxy)-4-oxo-3-[2-(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetylamino]-pentanoic acid;  
5-(2-Benzyl-3-phenyl-propionyloxy)-4-oxo-3-[2-(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetylamino]-pentanoic acid;  
4-Oxo-5-[(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetoxy]-3-[2-(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetylamino]-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-5-(3-benzyl-4-phenyl-butyryloxy)-4-oxo-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-5-(4-benzyl-5-phenyl-pentanoyloxy)-4-oxo-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-4-oxo-5-[(1-oxo-1,2,3,4-tetrahydro-naphthalen-2-yl)-acetoxy]-pentanoic acid; and  
5-(3-Benzyl-4-phenyl-butyryloxy)-3-[3-methyl-2-(3-phenyl-propionylamino)-butyrylamino]-4-oxo-pentanoic acid.--

*04  
Cancelled*

Please cancel claim 36, without waiver or prejudice, and insert therefor new claim 47:

--47. The compounds:

3-[2-(2-Benzylloxycarbonylamino-4-carboxy-butyrylamino)-3-methyl-butyrylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;  
3-[2-(2-Benzylloxycarbonylamino-3-methyl-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;  
3-(2-Acetylamino-3-methyl-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;  
3-[2-(2-Benzylloxycarbonylamino-3-methyl-butyrylamino)-propionylamino]-5-(3,3-diphenyl-propionyloxy)-4-oxo-pentanoic acid;  
3-[2-(2-Benzylloxycarbonylamino-3-methyl-butyrylamino)-propionylamino]-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid;  
3-[2-(2-Benzylloxycarbonylamino-3-methyl-butyrylamino)-propionylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

*05*

5-(2-Benzyl-3-phenyl-propionyloxy)-3-{2-[4-carboxy-2-(3-phenyl-propionylamino)-butyrylamino]-3-methyl-butyrylamino}-4-oxo-pentanoic acid;

3-(2-Benzylloxycarbonylamino-3-methyl-butyrylamino)-5-(3,3-diphenyl-propionyloxy)-4-oxo-pentanoic acid;

3-(2-Acetylamino-3-hydroxy-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Acetylamino-3-hydroxy-butyrylamino)-5-(3,3-diphenyl-propionyloxy)-4-oxo-pentanoic acid; and

5-(3,3-Diphenyl-propionyloxy)-4-oxo-3-[2-(4-phenyl-butyrylamino)-propionylamino]-pentanoic acid.--

*Please cancel* claim 37, without waiver or prejudice, and insert therefor new claim 48:

--48. The compound:

3-(2-{2-Acetylamino-3-(4-hydroxy-phenyl)-propionylamino}-4-carboxy-butyrylamino)-3-methyl-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid.--

*Please cancel* claim 42, without waiver or prejudice, and insert therefor new claim 49:

--49. The compounds:

3-(3-Carbamoyl-2-methyl-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Benzylloxycarbonylamino-3-methyl-naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-[(2-Carbamoyl-cyclopentanecarbonyl)-amino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-{2-Acetylamino-3-(4-hydroxy-phenyl)-propionylamino}-4-carboxy-butyrylamino)-3-methyl-butyrylamino)-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid;

3-(3-Carbamoyl-2-methyl-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(2-Carbamoylmethyl-3-methyl-butyrylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

3-(3-Benzyl-2-ureido-propionylamino)-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid;

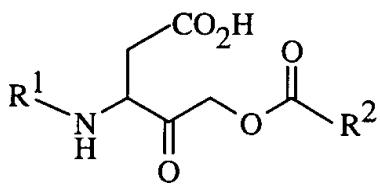
3-[2-(2-Benzylloxycarbonylamino-4-carboxy-butyrylamino)-3-methyl-butyrylamino]-5-(2-benzyl-3-phenyl-propionyloxy)-4-oxo-pentanoic acid;

*1*  
*Concl'd*

3-[2-[4-Carboxy-2-(3-phenyl-propionylamino)-butyrylamino]-3-methylbutyrylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid; and  
3-[2-(2-Acetylamino-4-carboxy-butyrylamino)-3-methyl-butyrylamino]-5-(naphthalen-1-yl-acetoxy)-4-oxo-pentanoic acid.--

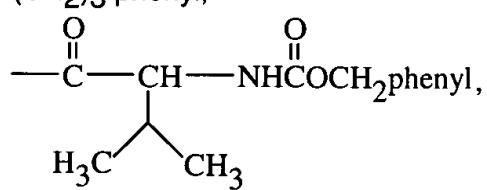
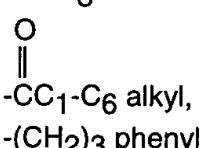
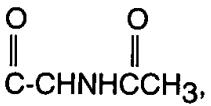
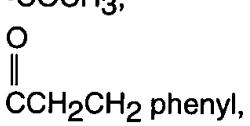
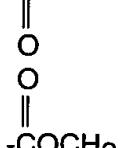
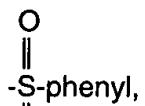
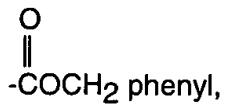
Please cancel claim 43, without waiver or prejudice, and insert therefor new claim 50:

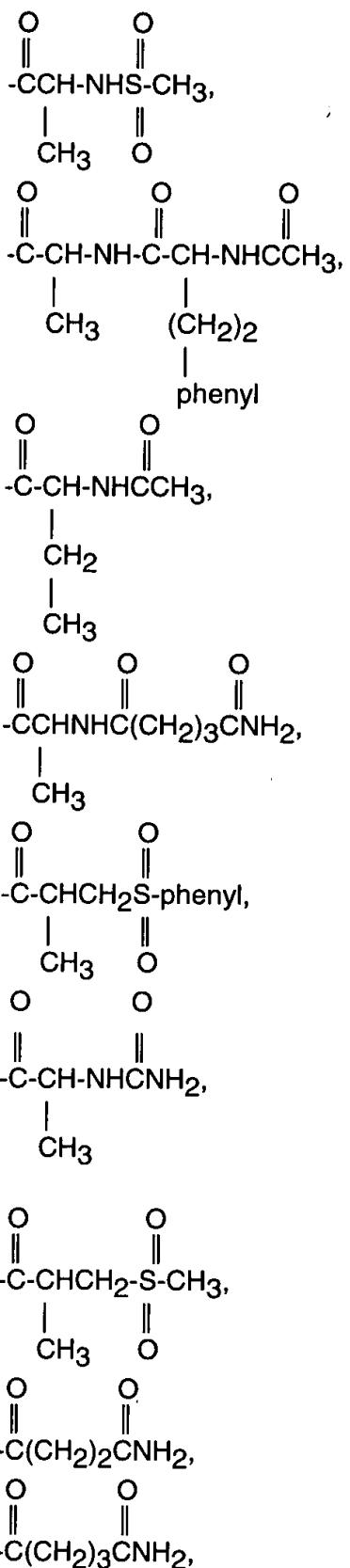
--50. A compound of the Formula I

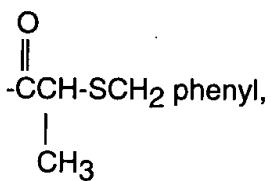
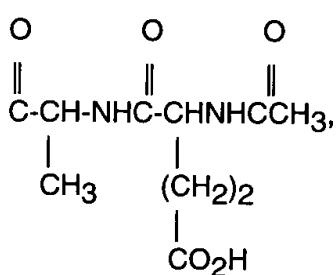
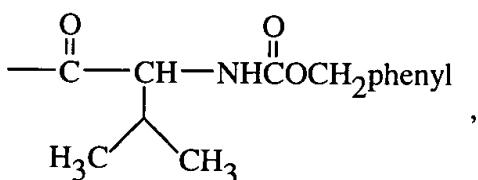
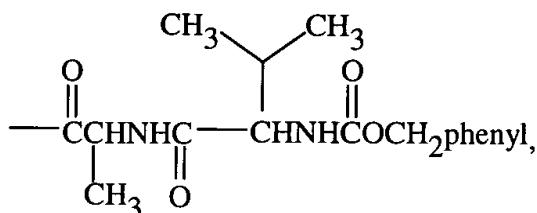
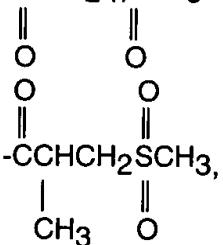
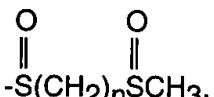
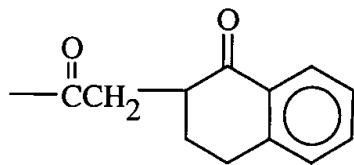
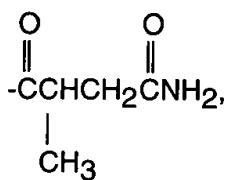


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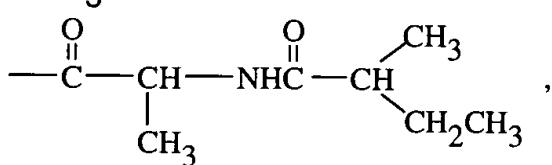
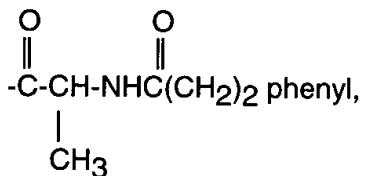
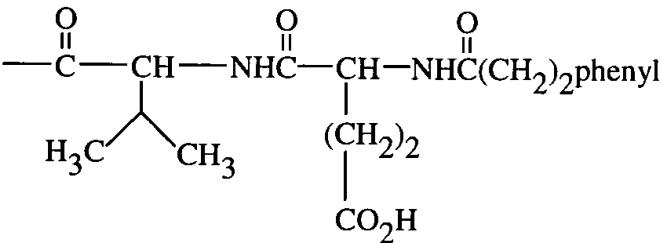
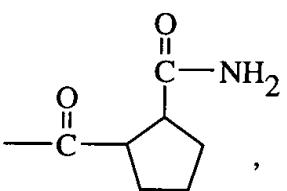
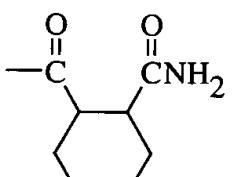
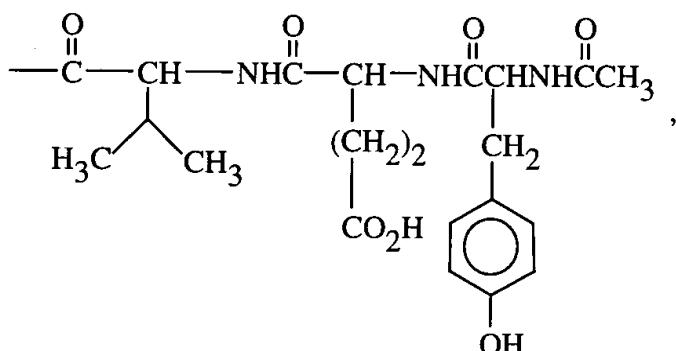
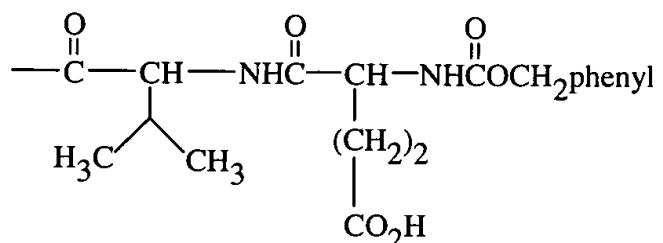
wherein  $\text{R}^1$  is



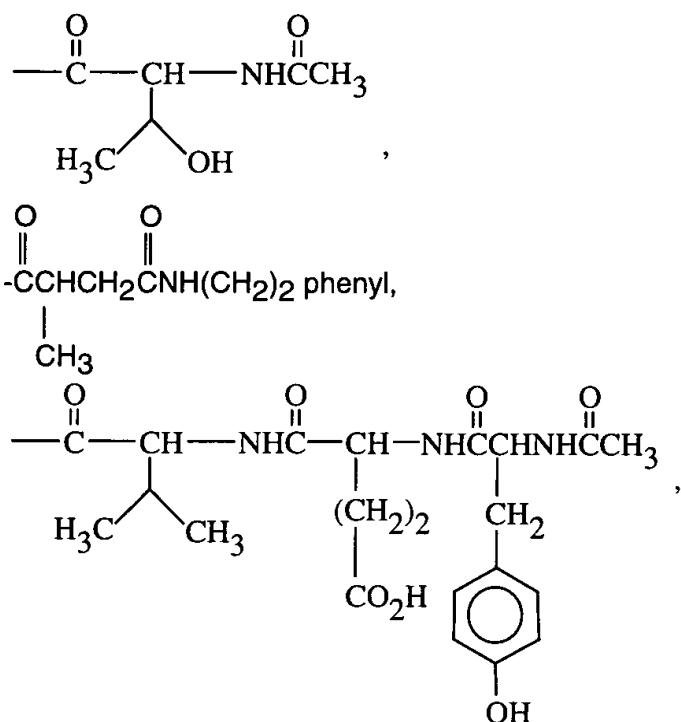




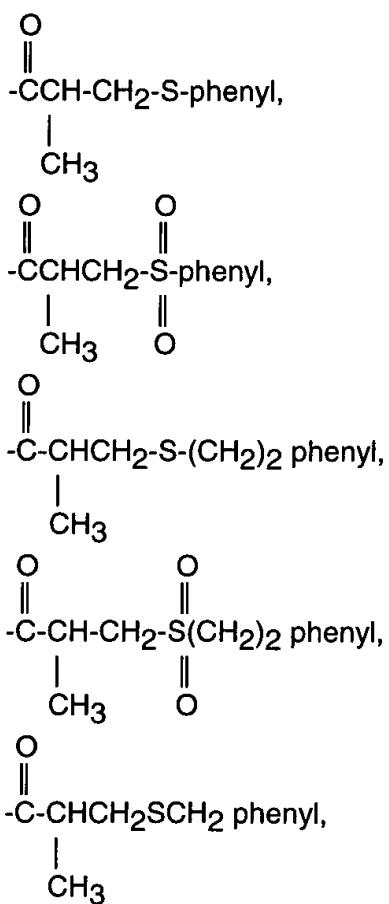
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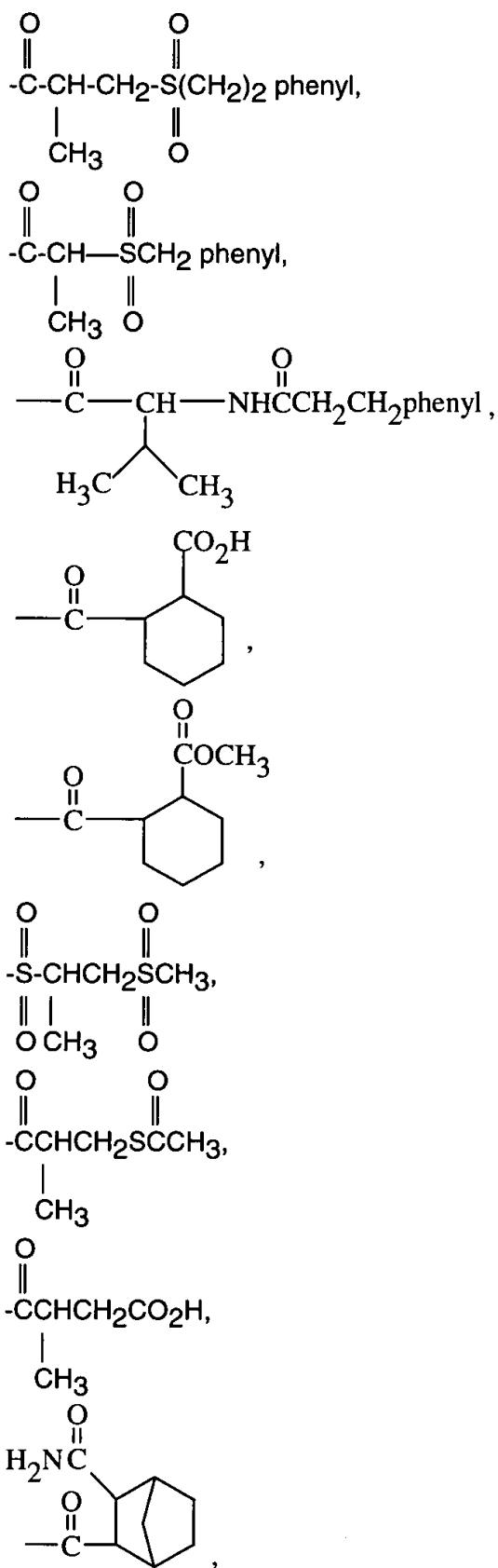


8  
O<sub>4</sub>H<sub>10</sub>

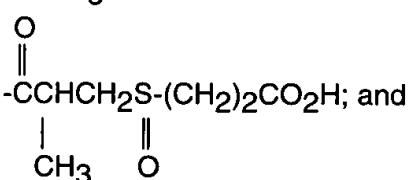
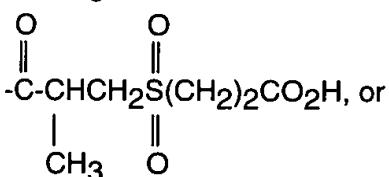
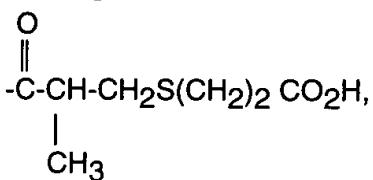
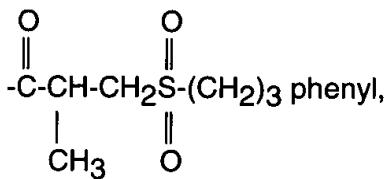
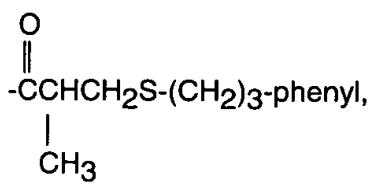


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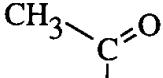




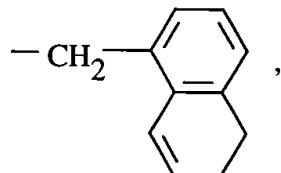
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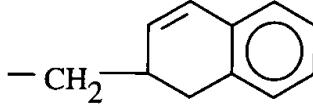


*08  
cont'd*  
R<sup>2</sup> is

- CH<sub>2</sub>CH<sub>2</sub> phenyl,
- CH<sub>2</sub> naphthyl,
- CH<sub>2</sub>CH<sub>2</sub> cyclohexyl,
- CH<sub>2</sub>O naphthyl,
- CH<sub>2</sub>O phenyl,
- CH<sub>2</sub>S-phenyl,
- CH<sub>2</sub>-substituted naphthyl,
- CH<sub>2</sub>CH(phenyl)<sub>2</sub>,
- (CH<sub>2</sub>)<sub>3</sub>-phenyl,
- CH-naphthyl,  
|  
CH<sub>3</sub>
- CH<sub>3</sub>   
— CH<sub>2</sub>—N— phenyl ,
- CH[CH<sub>2</sub>phenyl]<sub>2</sub>,

-CH-naphthyl,  
|  
OH  
-CH<sub>2</sub>-NH phenyl,  
— CH<sub>2</sub>—CH  
|      CH<sub>2</sub> substituted phenyl  
|      naphthyl  
-CH<sub>2</sub>-naphthyl-phenyl,  
-CH<sub>2</sub>-fluorenyl,  
-CH<sub>2</sub>-naphthyl-CH<sub>2</sub> phenyl,  
-CH<sub>2</sub>-substituted phenyl,



— CH<sub>2</sub>—CH  
|      phenyl  
|      CH<sub>2</sub>phenyl  
— CH<sub>2</sub>—CH  
|      phenyl  
|      NHphenyl  
— CH<sub>2</sub>—  
— CH<sub>2</sub>—